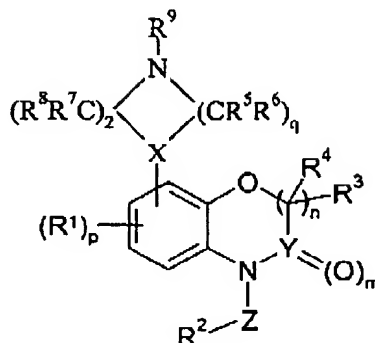


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Claim Listing

1. (Previously Presented) A compound of the formula:



or a pharmaceutically acceptable salt or prodrug thereof,
 wherein:

Y is C;

m is 1;

n is 1;

p is from 0 to 3;

q is from 1 to 3;

Z is $-(CR^aR^b)_r-$ or $-SO_2-$, where each of R^a and R^b is independently

hydrogen or alkyl;

r is from 0 to 2;

X is CH or N;

each R^1 is independently halo, alkyl, haloalkyl, heteroalkyl, alkoxy, cyano, $-S(O)_s-R^c$, $-C(=O)-NR^cR^d$, $-SO_2-NR^cR^d$, $-N(R^c)-C(=O)-R^d$, or $-C(=O)R^e$, where each of R^c and R^d is independently hydrogen or alkyl;

s is from 0 to 2;

R^2 is aryl or heteroaryl;

each of R^3 and R^4 is independently hydrogen, alkyl, hydroxyalkyl or alkoxyalkyl, or R^3 and R^4 together with their shared carbon may form a carbocyclic ring of 3 to 6 members; and

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each of R^5 , R^6 , R^7 , R^8 and R^9 is independently hydrogen or alkyl, or one of R^5 and R^6 together with one of R^7 , R^8 and R^9 and the atoms therebetween may form a ring of 5 to 7 members.

2. (Original) The compound of claim 1, wherein Z is $-(CR^aR^b)_r$.
3. (Original) The compound of claim 2, wherein X is N and q is 2.
4. (Canceled)
5. (Previously Presented) The compound of claim 3, wherein r is 1.
6. (Original) The compound of claim 5, wherein R^a and R^b are hydrogen.
7. (Original) The compound of claim 6, wherein R^2 is optionally substituted phenyl or optionally substituted naphthyl.
8. (Original) The compound of claim 7, wherein R^2 is 2-halophenyl, 3-halophenyl, 4-halophenyl, naphthyl-2-yl, 3-cyanophenyl, 4-cyanophenyl, 3-nitrophenyl, 3-aminophenyl, 3-methoxyphenyl, 3-ureaphenyl, or 3-methylsulfonylamino-phenyl.
9. (Original) The compound of claim 7, wherein p is 1 and R^1 is halo, methyl or methoxy.
10. (Original) The compound of claim 7, wherein R^3 and R^4 are hydrogen.
11. (Original) The compound of claim 7, wherein R^3 and R^4 are methyl.
12. (Original) The compound of claim 7, wherein one of R^3 and R^4 is hydrogen and the other is methyl.

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13. (Currently Amended) The compound of claim 7, wherein R³ and R⁴ together with the carbon atom therebetween form a cyclobutyl,

14. (Previously Presented) The compound of claim 8, wherein said compound is selected from:

4-benzyl-6-methyl-8-piprazin-1-yl-*H*-benzo[1,4]oxazin-3-one;
4-benzyl-6-methoxy-8-piperazin-1-yl-*H*-benzo[1,4]oxazin-3-one;
4-(2-fluoro-benzyl)-6-methoxy-8-piperazin-1-yl-4*H*-benzo[1,4]oxazin-3-one;
4-(2-chloro-benzyl)-6-methoxy-8-piperazin-1-yl-4*H*-benzo[1,4]oxazin-3-one;
4-(3-chloro-benzyl)-6-methoxy-8-piperazin-1-yl-4*H*-benzo[1,4]oxazin-3-one;
4-benzyl-8-piperazin-1-yl-4*H*-benzo[1,4]oxazin-3-one;
4-benzyl-6-fluoro-8-piperazin-1-yl-4*H*-benzo[1,4]oxazin-3-one;
4-(2-fluoro-benzyl)-8-piprazin-1-yl-4*H*-benzo[1,4]oxazin-3-one;
4-(4-fluoro-benzyl)-8-piperazin-1-yl-4*H*-benzo[1,4]oxazin-3-one;
4-(4-chloro-benzyl)-8-piperazin-1-yl-4*H*-benzo[1,4]oxazin-3-one;
4-(4-fluoro-benzyl)-6-fluoro-8-piperazin-1-yl-4*H*-benzo[1,4]oxazin-3-one;
4-(2-fluoro-benzyl)-6-fluoro-8-piperazin-1-yl-4*H*-benzo[1,4]oxazin-3-one;
4-(2-chloro-benzyl)-6-fluoro-8-piperazin-1-yl-4*H*-benzo[1,4]oxazin-3-one;
4-(4-chloro-benzyl)-6-fluoro-8-piprazin-1-yl-4*H*-benzo[1,4]oxazin-3-one;
6-fluoro-4-naphthalen-2-ylmethyl-8-piperazin-1-yl-4*H*-benzo[1,4]oxazin-3-one;
4-(3-chloro-benzyl)-6-fluoro-8-piperazin-1-yl-4*H*-benzo[1,4]oxazin-3-one;
3-(3-oxo-8-piperazin-1-yl-2,3-dihydro-benzo[1,4]oxazin-4-ylmethyl)-benzonitrile;
4-(3-fluoro-benzyl)-8-piprazin-1-yl-4*H*-benzo[1,4]oxazin-3-one;
4-benzyl-2,2-dimethyl-8-piperazin-1-yl-4*H*-benzo[1,4]oxazin-3-one;
(*R*)-4-benzyl-2-methyl-8-piprazin-1-yl-4*H*-benzo[1,4]oxazin-3-one;
4-benzyl-6-piprazin-1-yl-4*H*-benzo[1,4]oxazin-3-one;
4-(4-Fluoro-benzyl)-6-piperazin-1-yl-4*H*-benzo[1,4]oxazin-3-one;
(*S*)-4-Benzyl-2-methyl-8-piperazin-1-yl-4*H*-benzo[1,4]oxazin-3-one;
8-Piperazin-1-yl-4-pyridin-4-ylmethyl-4*H*-benzo[1,4]oxazin-3-one;
4-Benzyl-6-methyl-8-(4-methyl-piperazin-1-yl)-4*H*-benzo[1,4]oxazin-3-one;

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4-Benzyl-8-(4-methyl-piperazin-1-yl)-4*H*-benzo[1,4]oxazin-3-one;
4-(1-Phenyl-ethyl)-8-piperazin-1-yl-4*H*-benzo[1,4]oxazin-3-one;
4-(3-Methoxy-benzyl)-8-piperazin-1-yl-4*H*-benzo[1,4]oxazin-3-one;
4-(3-Nitro-benzyl)-8-piperazin-1-yl-4*H*-benzo[1,4]oxazin-3-one;
4-(3-Amino-benzyl)-8-piperazin-1-yl-4*H*-benzo[1,4]oxazin-3-one;
3-(3-Oxo-8-piperazin-1-yl-2,3-dihydro-benzo[1,4]oxazin-4-ylmethyl)-benzonitrile;
N-[3-(3-Oxo-8-piperazin-1-yl-2,3-dihydro-benzo[1,4]oxazin-4-ylmethyl)-phenyl]-methanesulfonamide;
4-(4-Fluoro-benzyl)-2,2-dimethyl-8-piperazin-1-yl-4*H*-benzo[1,4]oxazin-3-one;
4-(3-Fluoro-benzyl)-2,2-dimethyl-8-piperazin-1-yl-4*H*-benzo[1,4]oxazin-3-one;
[3-(3-Oxo-8-piperazin-1-yl-2,3-dihydro-benzo[1,4]oxazin-4-ylmethyl)-phenyl]-urea;
4-(3-Chloro-benzyl)-2,2-dimethyl-8-piperazin-1-yl-4*H*-benzo[1,4]oxazin-3-one;
4-Benzyl-8-(3,5-dimethyl-piperazin-1-yl)-4*H*-benzo[1,4]oxazin-3-one;
4-(4-Chloro-benzyl)-2,2-dimethyl-8-piperazin-1-yl-4*H*-benzo[1,4]oxazin-3-one;
4-Benzyl-6-fluoro-2,2-dimethyl-8-piperazin-1-yl-4*H*-benzo[1,4]oxazin-3-one;
4-(4-Chloro-benzyl)-6-fluoro-2,2-dimethyl-8-piperazin-1-yl-4*H*-benzo[1,4]oxazin-3-one;
6-Fluoro-4-(3-fluoro-benzyl)-2,2-dimethyl-8-piperazin-1-yl-4*H*-benzo[1,4]oxazin-3-one;
6-Fluoro-4-(2-fluoro-benzyl)-2,2-dimethyl-8-piperazin-1-yl-4*H*-benzo[1,4]oxazin-3-one;
6-Fluoro-4-(4-fluoro-benzyl)-2,2-dimethyl-8-piperazin-1-yl-4*H*-benzo[1,4]oxazin-3-one;
4-(3-Chloro-benzyl)-6-fluoro-2,2-dimethyl-8-piperazin-1-yl-4*H*-benzo[1,4]oxazin-3-one;
4-Benzyl-8-(3,3-dimethyl-piperazin-1-yl)-4*H*-benzo[1,4]oxazin-3-one;
4-Benzyl-2,2-spiro-cyclobutan-8-piperazin-1-yl-4*H*-benzo[1,4]oxazin-3-one.

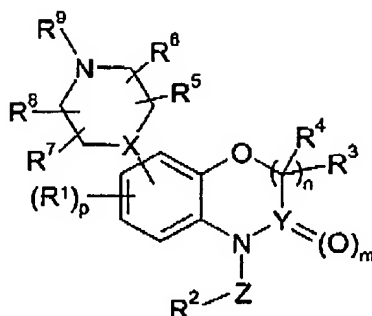
15. (Original) The compound of claim 6, wherein R² is heteroaryl.

16. (Original) The compound of claim 15, wherein R² is pyridine-4-yl.

17-32. (Canceled).

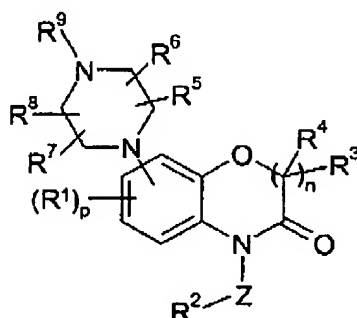
33. (Original) The compound of claim 1, wherein said compound is of the formula:

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or a pharmaceutically acceptable salt or prodrug thereof, wherein X, Y, Z, R¹, R², R³, R⁴, R⁵, R⁶, R⁷, R⁸, R⁹, m, n, and p are as defined in claim 1.

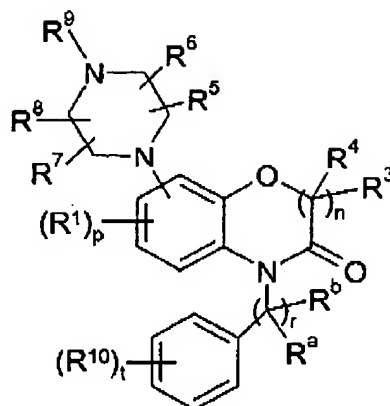
34. (Original) The compound of claim 1, wherein said compound is of the formula:



or a pharmaceutically acceptable salt or prodrug thereof, wherein Z, R¹, R², R³, R⁴, R⁵, R⁶, R⁷, R⁸, R⁹, n, and p are as defined in claim 1.

35. (Previously Presented) The compound of claim 1, wherein said compound is of the formula:

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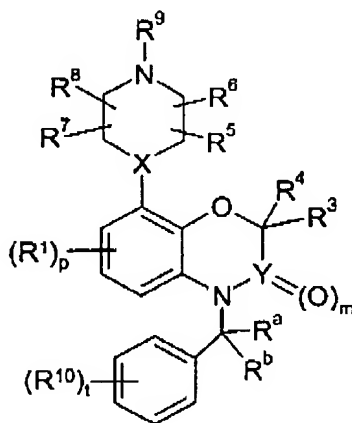


or a pharmaceutically acceptable salt or prodrug thereof, wherein R^1 , R^3 , R^4 , R^5 , R^6 , R^7 , R^8 , R^9 , R^a , R^b , n , p and r are as defined in claim 1, and wherein:

t is from 0 to 4; and

each R^{10} independently is halo, alkyl, alkoxy or cyano.

36. (Previously Presented) The compound of claim 1, wherein said compound is of the formula:



wherein X , Y , R^1 , R^3 , R^4 , R^5 , R^6 , R^7 , R^8 , R^9 , R^a , R^b , m , p and t are as recited in claim 1, and wherein:

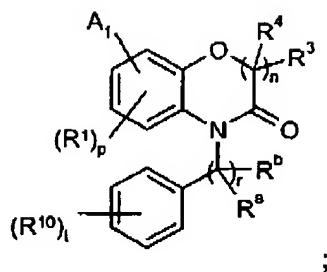
t is from 0 to 4; and

each R^{10} independently is halo, alkyl, alkoxy or cyano.

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37. (Original) The compound of claim 36, wherein R^1 is halo, methyl or methoxy.
38. (Original) The compound of claim 36 wherein R^3 and R^4 each independently is hydrogen or methyl.
39. (Original) The compound of claim 36, wherein R^3 and R^4 together with their shared carbon form a cyclobutyl group.
40. (Original) The compound of claim 36, wherein R^6 , R^7 , R^8 , R^9 each independently is hydrogen or methyl.
41. (Original) The compound of claim 36, wherein R^a and R^b each independently is hydrogen or methyl.
42. (Original) The compound of claim 36, wherein each R^{10} is hydrogen, halo, nitro, cyano, amino, urea, methoxy or methanesulfonylamino.
43. (Original) A pharmaceutical composition comprising an efficacious amount of the compound of claim 1 in admixture with a pharmaceutically acceptable carrier.
44. (Canceled)
45. (Canceled)
46. (Canceled)
47. (Previously Presented) A method for producing a substituted benzoxazinone compound, said method comprising:
- (a) contacting an N-arylalkyl benzoxazinone of the formula:

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wherein:

A_1 is a leaving group,

n is 1;

p is from 0 to 3;

r is from 0 to 2;

t is from 0 to 4;

each of R^a and R^b is independently hydrogen or alkyl;

each R^1 is independently halo, alkyl, haloalkyl, heteroalkyl, alkoxy, cyano,

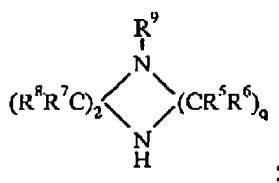
$-S(O)_s-R^c$, $-C(=O)-NR^cR^d$, $-SO_2-NR^cR^d$, $-N(R^c)-C(=O)-R^d$, or $-C(=O)$

R^c , where each of R^c and R^d is independently hydrogen or alkyl and s is from 0 to 2;

each of R^3 and R^4 is independently hydrogen or alkyl; and

each R^{10} is independently halo, alkyl, alkoxy or cyano;

with a heterocyclic compound of the formula:



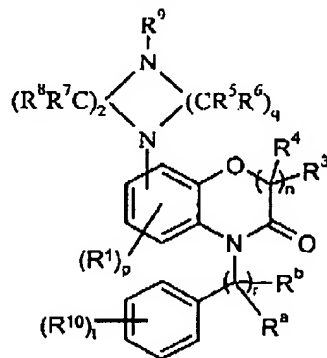
wherein:

q is from 1 to 3; and

each of R^5 , R^6 , R^7 , R^8 and R^9 is independently hydrogen or alkyl, or one of R^5 and R^6 together with one of R^7 , R^8 and R^9 may form a ring of 5 to 7 members;

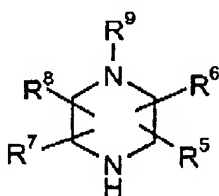
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in the presence of a palladium catalyst to produce the heterocyclyl-substituted N-arylalkyl benzoxaninone compound of the formula:

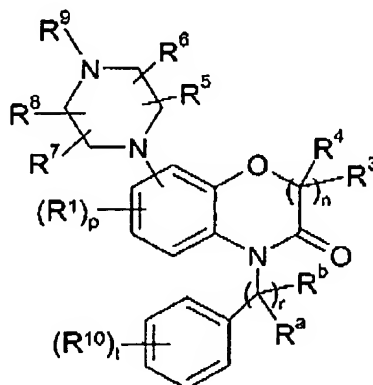


48. (Original) The method of claim 47, wherein the leaving groups A¹ is halo.

49. (Previously Presented) The method of claim 47, wherein the heterocyclic compound is of the formula:



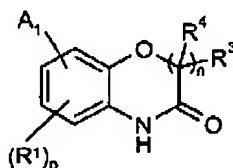
such that the heterocyclyl-substituted N-arylalkyl benzoxaninone compound is of the formula:



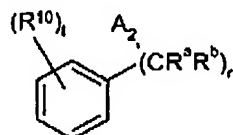
and R¹, R², R³, R⁴, R⁵, R⁶, R⁷, R⁸, R⁹, R¹⁰, n, p, r and t are as described in claim 47.

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50. (Original) The method of claim 47, further comprising:
 (a) contacting a benzoxazinone of the formula:



wherein n , p , A_1 , R^1 , R^3 and R^4 are as described in claim 1,
 with an alkylating agent of the formula:

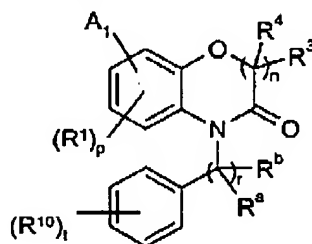


wherein:

A_2 is a leaving group and may be the same or different from A_1 ; and

r , t , R^a , R^b and R^{10} are as described in claim 41;

to produce the N-arylalkyl benzoxazinone of the formula:



51. (New) A method for enhancing cognitive memory in a subject, said method comprising administering to said subject a therapeutically effective amount of a compound of claim 1.